## **AMENDMENTS TO THE CLAIMS**

This listing of the claims will replace all prior versions, and listings, of claims in the application:

## **LISTING OF THE CLAIMS**

- 1. (currently amended) A compound having the structural formula  $L^1[MQ^1Q^2]L^2$  in which M is a mid-transition metal selected from the group consisting of Nb, Ta, Mo, W, Mn and Re,  $Q^1$  and  $Q^2$  are each a univalent radical substituent, and  $L^1$  and  $L^2$  are ligands coordinated to M, wherein each of  $L^1$  and  $L^2$  contains a first coordinating atom that is a nitrogen atom contained within a C=N group, and a second coordinating atom-that is either a second nitrogen-atom, optionally present-in a-second-C=N-group, or an oxygen, sulfur or phosphorus atom.
- 2. (original) The compound of claim 1, wherein, in each of  $L^1$  and  $L^2$ , the second coordinating atom is a second nitrogen atom.
- 3. (original) The compound of claim 2, wherein, in each of  $L^1$  and  $L^2$ , the second nitrogen atom is present in a second C=N group.
  - 4. (original) The compound of claim 3, wherein  $L^1$  and  $L^2$  are identical.
- 5. (original) The compound of claim 4, wherein the first nitrogen atom in each of  $L^1$  and  $L^2$  is bound to a first substituent  $R_S$ , and the second nitrogen atom in each of  $L^1$  and  $L^2$  is bound to a second substituent  $R_L$ , wherein the difference in steric bulk between  $R_S$  and  $R_L$  is sufficient to result in isospecificity when the compound is used as a polymerization catalyst.
  - 6. (original) The compound of claim 3, wherein  $L^1$  and  $L^2$  are different.
- 7. (original) The compound of claim 6, wherein the first and second nitrogen atoms in the ligand  $L^1$  are bound to a first substituent  $R_S$ , and the first and second nitrogen atoms in the ligand  $L^2$  are bound to a second substituent  $R_L$ , wherein the difference in steric bulk between  $R_S$  and  $R_L$  is sufficient to result in syndiospecificity when the compound is used as a polymerization catalyst.

- **8.** (original) The compound of claim 1, wherein the compound has a positive charge +y and is associated with y/z anions each bearing a negative charge -z.
- 9. (original) The compound of claim 8, wherein y and z are independently integers in the range of 1 to 4 inclusive.
  - 10. (original) The compound of claim 9, wherein y and z are independently 1 or 2.
- 11. (original) The compound of claim 8, wherein the anions are selected from the group consisting of halide and pseudohalide.
  - 12. (currently amended) A compound having the structure of formula (I)

(I)
$$\begin{bmatrix}
R^1 \\
R^2 \\
(CR^3R^4)_m
\end{bmatrix}$$

$$\begin{bmatrix}
Q^1 \\
Q^1
\end{bmatrix}$$

$$\begin{bmatrix}
Q^1 \\
Q^2
\end{bmatrix}$$

$$\begin{bmatrix}
L^1 \\
\end{bmatrix}$$

wherein:

M is a mid-transition metal selected from Groups VA, VIA and VIIA of the periodic table of the elements;

 $Q^1$  and  $Q^2$  are independently selected from the group consisting of hydrido, halide, alkoxy, amido, unsubstituted  $C_1$ - $C_{30}$  hydrocarbyl,  $C_1$ - $C_{30}$  hydrocarbyl substituted with one or more substituents such as electron withdrawing groups, and  $C_1$ - $C_{30}$  hydrocarbyl-substituted Group IVB elements, or  $Q^1$  and  $Q^2$  may together form an alkylidene olefin, acetylene, or a five- or six-membered cyclic hydrocarbyl group;

m and n are independently zero or 1; q is an optional double bond; X is N, O, S or P, with the provisos that (a) when X is N or P, then either n is 1 or q is present as a double bond, but not both, and (b) when X is O or S, then n is zero and q is absent;

R<sup>1</sup>, R<sup>6</sup>, and R<sup>7</sup> are independently hydrido, hydrocarbyl or substituted hydrocarbyl, and R<sup>2</sup> and R<sup>5</sup> are independently hydrido, halo, hydrocarbyl or substituted hydrocarbyl, or R<sup>1</sup> and R<sup>2</sup> and/or R<sup>5</sup> and R<sup>6</sup> may be taken together to form a linkage -Q-, resulting in a five- or six-membered ring, wherein Q is -[(CR)<sub>a</sub>(Z)<sub>b</sub>]- in which a is 2, 3 or 4, Z is N, O or S, b is zero or 1, the sum of a and b is 3 or 4, and R is selected from the group consisting of hydrido, halo, hydrocarbyl, hydrocarbyloxy, trialkylsilyl, NR<sup>8</sup><sub>2</sub>, OR<sup>9</sup>, and NO<sub>2</sub>, wherein R<sup>8</sup> and R<sup>9</sup> are each independently hydrocarbyl, or wherein R moieties on adjacent carbon atoms may be linked to form an additional five- or six-membered ring, or R<sup>2</sup> and R<sup>5</sup> may together form a-linkage -Q--as-just defined;

 $R^3$  and  $R^4$  are independently selected from the group consisting of hydrido and hydrocarbyl, or at least one of  $R^3$  and  $R^4$  may be bound through a lower alkylene linkage to an atom contained within  $L^A$  or  $L^B$ ;

 $L^A$  and  $L^B$  are ligands which may be the same or different and are independently selected from the group consisting of nitrogen-containing, sulfur-containing and oxygen-containing heterocycles, tertiary amines and phosphines, or  $L^A$  and  $L^B$  may together form a single bidentate ligand that may or may not be the same as  $L^1$ ,

with the proviso that when (a)  $L^A$  and  $L^B$  form a single bidentate ligand that is identical to  $L^1$  and M is V or Cr, then either (b)  $R^1$  and  $R^2$  or  $R^5$  and  $R^6$  are taken together to form a linkage -Q- as defined above, or (c) X is other than N, or both (b) and (c).

- 13. (original) The compound of claim 12, wherein the compound has a positive charge +y and is associated with y/z anions each bearing a negative charge -z.
- 14. (original) The compound of claim 13, wherein y and z are independently integers in the range of 1 to 4 inclusive.
  - 15. (original) The compound of claim 14, wherein y and z are independently 1 or 2.
- 16. (currently amended) The compound of claim 12-13, wherein the anions are selected from the group consisting of halide and pseudohalide.

## 17. (currently amended) The compound of claim 12, having the structure of formula (II)

wherein,  $q^a$ , ma, na, and  $R^{1a}$  through  $R^{7a}$  are defined as for q, m, n and  $R^1$  through  $R^7$ , respectively.

18. (original) The compound of claim 17, wherein the compound has a positive charge +y and is associated with y/z anions each bearing a negative charge -z.

- 19. (original) The compound of claim 18, wherein y and z are independently integers in the range of 1 to 4 inclusive.
  - 20. (original) The compound of claim 19, wherein y and z are independently 1 or 2.
- 21. (currently amended) The compound of claim 17-18, wherein the anions are selected from the group consisting of halide and pseudohalide.
  - 22. (original) The compound of claim 17, having the structure of formula (V)

$$(V) \qquad \begin{array}{c} R^{21} \\ R^{20} \\ R^{20a} \\ R^{20a} \\ R^{21a} \\ R^{21a} \\ R^{22a} \\ R^{23a} \\ R^{5a} \\ R^{5a} \end{array}$$

wherein:

 $R^{20}$ ,  $R^{20a}$ ,  $R^{21}$ ,  $R^{21a}$ ,  $R^{22}$ ,  $R^{22a}$ ,  $R^{23}$  and  $R^{23a}$  are hydrido or hydrocarbyl of 1 to 10 carbon atoms, or any two adjacent  $R^{20}$ ,  $R^{20a}$ ,  $R^{21}$ ,  $R^{21a}$ ,  $R^{22}$ ,  $R^{22a}$ ,  $R^{23}$  and  $R^{23a}$  groups may be linked to form a five- or six-membered aromatic ring.

- 23. (original) The compound of claim 22, wherein  $R^{20}$ ,  $R^{20a}$ ,  $R^{21}$ ,  $R^{21a}$ ,  $R^{22}$ ,  $R^{22a}$ ,  $R^{23}$  and  $R^{23a}$  are hydrido.
- **24.** (original) The compound of claim 22, wherein  $R^{20}$  and  $R^{20a}$  are methyl, and  $R^{21}$ ,  $R^{21a}$ ,  $R^{22}$ ,  $R^{22a}$ ,  $R^{23}$  and  $R^{23a}$  are hydrido.
- **25.** (original) The compound of any one of claims 21, 22 or 23, wherein the compound has a positive charge +y and is associated with y/z anions each bearing a negative charge -z.

- 26. (original) The compound of claim 25, wherein y and z are independently integers in the range of 1 to 4 inclusive.
  - 27. (original) The compound of claim 26, wherein y and z are independently 1 or 2.
- 28. (original) The compound of claim 25, wherein the anions are selected from the group consisting of halide and pseudohalide.
  - 29. (currently amended) A compound having the structure of formula (III)

(III) 
$$\begin{bmatrix} R^{12} \\ R^{13} \end{bmatrix}_{j} = \begin{bmatrix} R^{10} \\ R^{11} \end{bmatrix}$$

wherein:

M is a mid-transition metal selected from Groups VA, VIA and VIIA of the periodic table of the elements;

 $Q^1$  and  $Q^2$  are independently selected from the group consisting of hydrido, halide, alkoxy, amido, unsubstituted  $C_1$ - $C_{30}$  hydrocarbyl,  $C_1$ - $C_{30}$  hydrocarbyl substituted with one or more substituents such as electron-withdrawing groups, and  $C_1$ - $C_{30}$  hydrocarbyl-substituted Group IVB elements, or  $Q^1$  and  $Q^2$  may together form an alkylidene olefin, acetylene, or a five- or six-membered cyclic hydrocarbyl group;

L<sup>A</sup> and L<sup>B</sup> are ligands which may be the same or different and are independently selected from the group consisting of nitrogen-containing, sulfur-containing and oxygen-containing heterocycles, tertiary amines and phosphines, or L<sup>A</sup> and L<sup>B</sup> may together form a single bidentate ligand that may or may not be the same as L<sup>3</sup>;

i and j are independently zero, 1, 2 or 3; and

- R<sup>10</sup>, R<sup>11</sup>, R<sup>12</sup> and R<sup>13</sup> are independently hydrido, hydrocarbyl or substituted hydrocarbyl.
- **30.** (withdrawn) A catalyst system comprising the compound of any one of claims 1, 5, 7, 12, 17 or 22 and a catalyst activator effective to produce a catalytically active ionic species when combined with said compound.
- 31. (withdrawn) The catalyst system of claim 30, wherein the catalyst activator is aluminum-containing or boron-containing.
- 32. (withdrawn) The catalyst-system of claim 31, wherein the catalyst activator is aluminum containing.
- **33.** (withdrawn) The catalyst system of claim 32, wherein the catalyst activator is an organoaluminum compound.
- 34. (withdrawn) The catalyst system of claim 33, wherein the catalyst activator is an alkyl aluminoxane.
- **35.** (withdrawn) The catalyst system of claim 34, wherein the catalyst activator is methyl aluminoxane.
- **36.** (withdrawn) The catalyst system of claim 31, wherein the catalyst activator is boron-containing.
- 37. (withdrawn) The catalyst system of claim 36, wherein the catalyst activator is a fluorohydrocarbylboron compound.
- **38.** (withdrawn) The catalyst system of claim 37, wherein the catalyst activator is a fluorinated phenylborate.
- **39.** (withdrawn) The catalyst system of claim 30, further including an inert polymerization diluent.

**40.** (withdrawn) The catalyst system of claim 39, wherein the diluent is a volatile hydrocarbon solvent.

## Claims 41-51 (canceled).

- **52.** (new) The compound of claim 12, wherein  $Q^1$  and  $Q^2$  are independently selected from the group consisting of hydrido, halide, alkoxy, amido, unsubstituted  $C_1$ - $C_{30}$  hydrocarbyl,  $C_1$ - $C_{30}$  hydrocarbyl substituted with one or more electron-withdrawing groups, and  $C_1$ - $C_{30}$  hydrocarbyl-substituted Group IVB elements, or  $Q^1$  and  $Q^2$  may together form an alkylidene olefin, acetylene, or a five- or six-membered cyclic hydrocarbyl group.
- **53.** (new) The compound of claim 29, wherein  $Q^1$  and  $Q^2$  are independently selected from the group consisting of hydrido, halide, alkoxy, amido, unsubstituted  $C_1$ - $C_{30}$  hydrocarbyl,  $C_1$ - $C_{30}$  hydrocarbyl substituted with one or more electron-withdrawing groups, and  $C_1$ - $C_{30}$  hydrocarbyl-substituted Group IVB elements, or  $Q^1$  and  $Q^2$  may together form an alkylidene olefin, acetylene, or a five- or six-membered cyclic hydrocarbyl group.
- 54. (new) The compound of claim 1, wherein one of  $L^1$  and  $L^2$  contains a second coordinating atom that is either a second nitrogen atom, optionally present in a second C=N group, or an oxygen, sulfur or phosphorus atom, and the other of  $L^1$  and  $L^2$  contains a second coordinating atom that is either a second nitrogen atom, optionally present in a second C=N group, or a sulfur or phosphorus atom.